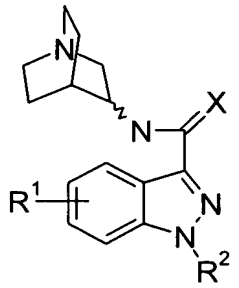
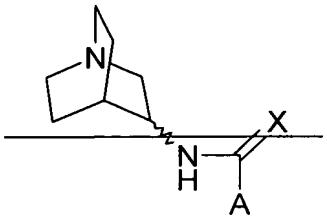
This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (Currently Amended): A compound of Formula I:



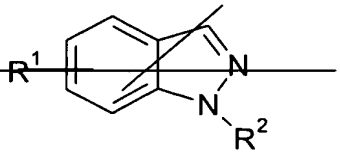
(I)



-(I)-

wherein

~~A is an indazolyl group according to subformula (a)~~

-(a) 

X is O or S;

R^1 is H, F, Cl, Br, I, OH, CN, nitro, NH_2, alkyl having 1 to 4 carbon atoms, fluorinated alkyl having 1 to 4 carbon atoms, cycloalkyl having 3 to 7 carbon atoms, cycloalkylalkyl having 4 to 7 carbon atoms, alkoxy having 1 to 4 carbon atoms, cycloalkoxy having 3 to 7 carbon atoms, cycloalkylalkoxy having 4 to 7 carbon atoms, alkylthio having 1 to 4 carbon atoms, fluorinated alkoxy having 1 to 4 carbon atoms, hydroxyalkyl ~~hydoxyalkyl~~ having 1 to 4 carbon atoms, hydroxyalkoxy having 2 to 4 carbon atoms, monoalkylamino having 1 to 4 carbon atoms, dialkylamino wherein each alkyl group independently has 1 to 4 carbon atoms, Ar or Het;

R^2 is H, alkyl having 1 to 4 carbon atoms, cycloalkyl having 3 to 7 carbon atoms, or cycloalkylalkyl having 4 to 7 carbon atoms;

Ar is an aryl group containing 6 to 10 carbon atoms which is unsubstituted or substituted one or more times by alkyl having 1 to 8 carbon atoms, alkoxy having 1 to 8 carbon atoms, halogen, dialkylamino wherein the alkyl portions each have 1 to 8 carbon atoms, amino, cyano, hydroxyl, nitro, halogenated alkyl having 1 to 8 carbon atoms, halogenated alkoxy having 1 to 8 carbon atoms, hydroxyalkyl having 1 to 8 carbon atoms, hydroxyalkoxy having 2 to 8 carbon atoms, alkenyloxy having 3 to 8 carbon atoms, alkylthio having 1 to 8 carbon atoms, alkylsulphinyl having 1 to 8 carbon atoms, alkylsulphonyl having 1 to 8 carbon atoms, monoalkylamino having 1 to 8 carbon atoms, cycloalkylamino wherein the cycloalkyl group has 3 to 7 carbon atoms and is optionally substituted, aryloxy wherein the aryl portion contains 6 to 10 carbon atoms and is optionally substituted, arylthio wherein the aryl portion contains 6 to 10 carbon atoms and is optionally substituted, cycloalkyloxy wherein the cycloalkyl group has 3 to 7 carbon atoms and is optionally substituted, sulfo, sulfonylamino, acylamido, acyloxy or combinations thereof; and

Het is a heterocyclic group, which is fully saturated, partially saturated or fully

unsaturated, having 5 to 10 ring atoms in which at least 1 ring atom is a N, O or S atom, which is unsubstituted or substituted one or more times by halogen, aryl having 6 to 10 carbon atoms and is optionally substituted, alkyl having 1 to 8 carbon atoms, alkoxy having 1 to 8 carbon atoms, cyano, trifluoromethyl, nitro, oxo, amino, monoalkylamino having 1 to 8 carbon atoms, dialkylamino wherein each alkyl group has 1 to 8 carbon atoms, or combinations thereof; or

a pharmaceutically acceptable salt thereof,

wherein if the compound exhibits chirality it can be in the form of a mixture of enantiomers such as a racemate or a mixture of diastereomers, or can be in the form of a single enantiomer or a single diastereomer~~, and~~

~~wherein the indazolyl group of group A is attached via its 3, 4, or 7 position~~.

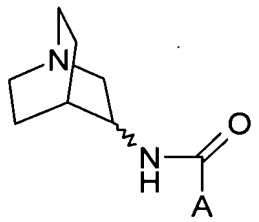
2. (Cancelled):

3. (Cancelled):

4. (Cancelled):

5. (Cancelled):

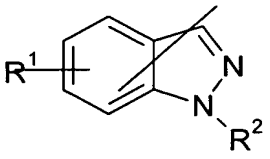
6. (Currently Amended): A compound according to Formula I':



(I')

wherein

A is an indazolyl according to subformula (a),

(a) 

R^1 is H, F, Cl, Br, I, OH, CN, nitro, NH_2, alkyl having 1 to 4 carbon atoms, fluorinated alkyl having 1 to 4 carbon atoms, cycloalkyl having 3 to 7 carbon atoms, cycloalkylalkyl having 4 to 7 carbon atoms, alkoxy having 1 to 4 carbon atoms, cycloalkoxy having 3 to 7 carbon atoms, alkylthio having 1 to 4 carbon atoms, fluorinated alkoxy having 1 to 4 carbon atoms, hydroxyalkyl ~~hydoxyalkyl~~ having 1 to 4 carbon atoms, hydroxyalkoxy having 2 to 4 carbon atoms, monoalkylamino having 1 to 4 carbon atoms, dialkylamino wherein each alkyl group independently has 1 to 4 carbon atoms, Ar or Het;

R^2 is H, alkyl having 1 to 4 carbon atoms, cycloalkyl having 3 to 7 carbon atoms, or cycloalkylalkyl having 4 to 7 carbon atoms;

Ar is an aryl group containing 6 to 10 carbon atoms which is unsubstituted or substituted one or more times by alkyl having 1 to 8 carbon atoms, alkoxy having 1 to 8 carbon atoms, halogen, dialkylamino wherein the alkyl portions each have 1 to 8 carbon atoms, amino, cyano, hydroxyl, nitro, halogenated alkyl having 1 to 8

carbon atoms, halogenated alkoxy having 1 to 8 carbon atoms, hydroxyalkyl having 1 to 8 carbon atoms, hydroxyalkoxy having 2 to 8 carbon atoms, alkenyloxy having 3 to 8 carbon atoms, alkylthio having 1 to 8 carbon atoms, alkylsulphinyl having 1 to 8 carbon atoms, alkylsulphonyl having 1 to 8 carbon atoms, monoalkylamino having 1 to 8 carbon atoms, cycloalkylamino wherein the cycloalkyl group has 3 to 7 carbon atoms and is optionally substituted, aryloxy wherein the aryl portion contains 6 to 10 carbon atoms and is optionally substituted, arylthio wherein the aryl portion contains 6 to 10 carbon atoms and is optionally substituted, cycloalkyloxy wherein the cycloalkyl group has 3 to 7 carbon ~~C~~ atoms and is optionally substituted, sulfo, sulfonylamino, acylamido, acyloxy or combinations thereof; and

Het is a heterocyclic group, which is fully saturated, partially saturated or fully unsaturated, having 5 to 10 ring atoms in which at least 1 ring atom is a N, O or S atom, which is unsubstituted or substituted one or more times by halogen, aryl having 6 to 10 carbon atoms and is optionally substituted, alkyl having 1 to 8 carbon atoms, alkoxy having 1 to 8 carbon atoms, cyano, trifluoromethyl, nitro, oxo, amino, monoalkylamino having 1 to 8 carbon atoms, dialkylamino wherein each alkyl group has 1 to 8 carbon atoms, or combinations thereof; or

a pharmaceutically acceptable salt thereof,

wherein if the compound exhibits chirality it can be in the form of a mixture of enantiomers such as a racemate or a mixture of diastereomers, or can be in the form of a single enantiomer or a single diastereomer, and

wherein the indazolyl group of subformula (a) is attached to the remainder of the compound via its 3~~, 4 or 7~~ position.

7. (Cancelled):

8. (Cancelled):

9. (Cancelled):

10. (Cancelled):

11. (Previously Presented): A compound according to claim 1, wherein R^1 is H, F, Cl, Br, 2-thiophenyl, 3-thiophenyl, 3-furyl, or phenyl.

12. (Previously Presented): A compound according to claim 1, wherein R^2 is H, methyl, 2-thiophenyl, 3-thiophenyl, 3-furyl, or phenyl.

13. (Cancelled):

14 (Previously Presented): A compound according to claim 1, wherein R^1 is H, F, Cl, Br, methyl, methoxy, or amino.

15. (Previously Presented): A compound according to claim 1, wherein R^2 is H or methyl.

16. (Cancelled):

17. (Cancelled):

18. (Cancelled):

19. (Previously Presented): A compound according to claim 1, wherein R^1 is H, F, Cl, Br, 2-thiophenyl, 3-thiophenyl, 3-furyl, or phenyl, and R^2 is H, methyl, 2-thiophenyl, 3-thiophenyl, 3-furyl, or phenyl.

20. (Cancelled):

21. (Currently Amended) A compound according to claim 1, wherein said compound is selected from:

N-(1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide,

N-((3R)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide,

1-Methyl-1H-Indazole-3-carboxamide, N-1-aza-bicyclo[2,2,2]oct-3-yl,

(R) 1-Methyl-1H-Indazole-3-carboxamide, N-1-aza-bicyclo[2,2,2]oct-3-yl,

(S) 1-Methyl-1H-Indazole-3-carboxamide, N-1-aza-bicyclo[2,2,2]oct-3-yl,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(bromo)-1H-indazole-3-carboxamide,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(cyclopropyl)-1H-indazole-3-carboxamide,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(bromo)-1H-indazole-3-carboxamide,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(bromo)-1H-indazole-3-carboxamide,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(furan-3-yl)-1H-indazole-3-carboxamide,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(phenyl)-1H-indazole-3-carboxamide,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-2-yl)-1H-indazole-3-carboxamide,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-3-yl)-1H-indazole-3-carboxamide,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(bromo)-1H-indazole-3-carboxamide,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(furan-3-yl)-1H-indazole-3-carboxamide,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(phenyl)-1H-indazole-3-carboxamide,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-2-yl)-1H-indazole-3-carboxamide,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-3-yl)-1H-indazole-3-carboxamide,
~~N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-4-carboxamide,~~
~~N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)- 1H-indazole-4-carboxamide,~~
~~N-(1H-Indazol-4-yl) 1-azabicyclo[2,2,2]oct-3-ylcarboxamide,~~
~~N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-7-carboxamide,~~
~~N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-7-carboxamide,~~
~~N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-7-carboxamide,~~
1-Aza-bicyclo[2,2,2]oct-3-ylcarboxamide, N-1H-indazol-3-yl,
(S) 1-Aza-bicyclo[2,2,2]oct-3-ylcarboxamide, N-1H-indazol-3-yl,
(R) 1-Aza-bicyclo[2,2,2]oct-3-ylcarboxamide, N-1H-indazol-3-yl,
~~(S) 1-Aza-bicyclo[2,2,2]oct-3-ylcarboxamide, N-1H-indazol-4-yl,~~
~~(R) 1-Aza-bicyclo[2,2,2]oct-3-ylcarboxamide, N-1H-indazol-4-yl,~~
~~1-Aza-bicyclo[2,2,2]oct-3-ylcarboxamide, N-1H-indazol-7-yl,~~
~~(S) 1-Aza-bicyclo[2,2,2]oct-3-ylcarboxamide, N-1H-indazol-7-yl,~~
~~(R) 1-Aza-bicyclo[2,2,2]oct-3-ylcarboxamide, N-1H-indazol-7-yl,~~
and pharmaceutically acceptable salts thereof.

22. (Previously Presented): A pharmaceutical composition comprising a compound according to claim 1 and a pharmaceutically acceptable carrier.

23. (Cancelled):

24. (Cancelled):

25. (Cancelled):

26. (Cancelled):

27. (Cancelled):

28. (Cancelled):

29. (Cancelled):

30. (Cancelled):

31. (Cancelled):

32. (Cancelled):

33. (Cancelled):

34. (Cancelled):

35. (Cancelled):

36. (Cancelled):

37. (Cancelled):

38. (Previously Presented): A compound according to claim 21, wherein said compound is in the form of a hydrochloride or hydroformate salt.

39. (Currently Amended): A compound according to claim 38, wherein said compound is selected from:

N-(-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide hydrochloride,

N-((3R)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide hydrochloride,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide hydrochloride,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(cyclopropyl)-1H-indazole-3-carboxamide hydroformate,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide hydroformate,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide hydroformate,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide hydroformate,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide hydroformate,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide hydroformate,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide hydroformate,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide hydroformate,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide hydroformate,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(furan-3-yl)-1H-indazole-3-carboxamide hydroformate,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(phenyl)-1H-indazole-3-carboxamide hydroformate,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-2-yl)-1H-indazole-3-carboxamide hydroformate,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-3-yl)-1H-indazole-3-carboxamide hydroformate,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(furan-3-yl)-1H-indazole-3-carboxamide hydroformate,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(phenyl)-1H-indazole-3-carboxamide hydroformate,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-2-yl)-1H-indazole-3-carboxamide hydroformate, and

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-3-yl)-1H-indazole-3-carboxamide

hydroformate~~,~~

~~N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-7-carboxamide hydrochloride, and~~

~~N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-7-carboxamide hydrochloride~~.

40. (Cancelled):

41. (Cancelled):

42. (Cancelled):

43. (Cancelled):

44. (Cancelled):

45. (Cancelled):

46. (Cancelled):

47. (Cancelled):

48. (Previously Presented): A compound according to claim 21, wherein said compound is N-(1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

49. (Previously Presented): A compound according to claim 48, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

50. (Previously Presented): A compound according to claim 48, wherein said

compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

51. (Previously Presented): A compound according to claim 21, wherein said compound is N-(1-Azabicyclo[2.2.2]oct-3-yl)-5-(bromo)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

52. (Previously Presented): A compound according to claim 51, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(bromo)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

53. (Previously Presented): A compound according to claim 51, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(bromo)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

54. (Previously Presented): A compound according to claim 21, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(cyclopropyl)-1H-indazole-3-carboxamide carboxamide or a pharmaceutically acceptable salt thereof.

55. (Previously Presented): A compound according to claim 21, wherein said compound is N-(1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

56. (Previously Presented): A compound according to claim 55, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

57. (Previously Presented): A compound according to claim 55, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide or

a pharmaceutically acceptable salt thereof.

58. (Previously Presented): A compound according to claim 21, wherein said compound is N-(1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

59. (Previously Presented): A compound according to claim 58, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

60. (Previously Presented): A compound according to claim 58, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

61. (Previously Presented): A compound according to claim 21, wherein said compound is N-(1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

62. (Previously Presented): A compound according to claim 61, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

63. (Previously Presented): A compound according to claim 61, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

64. (Previously Presented): A compound according to claim 21, wherein said compound is N-(1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

65. (Previously Presented): A compound according to claim 64, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

66. (Previously Presented): A compound according to claim 64, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

67. (New): A compound according to claim 39, wherein said compound is N-(-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide hydrochloride.

68. (New): A compound according to claim 39, wherein said compound is N-((3R)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide hydrochloride.

69. (New): A compound according to claim 39, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide hydrochloride.

70. (New): A compound according to claim 39, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(cyclopropyl)-1H-indazole-3-carboxamide hydroformate.

71. (New): A compound according to claim 39, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide hydroformate.

72. (New): A compound according to claim 39, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide hydroformate.

73. (New): A compound according to claim 39, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide hydroformate.

74. (New): A compound according to claim 39, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide hydroformate.

75. (New): A compound according to claim 39, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide hydroformate.

76. (New): A compound according to claim 39, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide hydroformate.

77. (New): A compound according to claim 39, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide hydroformate.

78. (New): A compound according to claim 39, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide hydroformate.

79. (New): A compound according to claim 11, wherein R^2 is H or methyl.

80. (New): A compound according to claim 1, wherein Ar is substituted or unsubstituted phenyl, naphthyl, or biphenyl, and Het is substituted or unsubstituted tetrahydrofuranyl, tetrahydrothienyl, pyrrolidinyl, piperidinyl, piperazinyl, morpholinyl, isoxazolinyl, furyl, thienyl, pyrrolyl, pyrazolyl, imidazolyl, pyridyl, pyrimidinyl, indolyl, quinolinyl, isoquinolinyl, or naphthyridinyl.